



Jupyter C3\_W2\_lecture\_notebook\_RNNs Last Checkpoint: a few seconds ago (autosaved)



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### Vanilla RNNs, GRUs and the scan function

In this notebook, you will learn how to define the forward method for vanilla RNNs and GRUs. Additionally, you will see how to define and use the function scan to compute forward propagation for RNNs.

By completing this notebook, you will:

- Be able to define the forward method for vanilla RNNs and GRUs
- Be able to define the scan function to perform forward propagation for RNNs
- · Understand how forward propagation is implemented for RNNs.

An implementation of the sigmoid function is provided below so you can use it in this notebook.

```
In [2]: H def sigmoid(x): # Sigmoid function
    return 1.0 / (1.0 + np.exp(-x))
```

## Part 1: Forward method for vanilla RNNs and GRUs

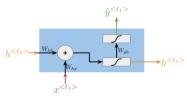
In this part of the notebook, you'll see the implementation of the forward method for a vanilla RNN and you'll implement that same method for a GRU. For this excersice you'll use a set of random weights and variables with the following dimensions:

- Embedding size ( emb ): 128
- Hidden state size ( h\_dim ): (16,1)

The weights  $w_a$  and biases  $b_a$  are initialized with dimensions ( $h_d$ im, emb +  $h_d$ im) and ( $h_d$ im, 1). We expect the hidden state  $h_t$  to be a column vector with size ( $h_d$ im, 1) and the initial hidden state  $h_t$ 0 is a vector of zeros.

### 1.1 Forward method for vanilla RNNs

The vanilla RNN cell is quite straight forward. Its most general structure is presented in the next figure:



As you saw in the lecture videos, the computations made in a vanilla RNN cell are equivalent to the following equations:

$$\begin{split} h^{< t>} &= g(W_h[h^{< t-1>}, x^{< t>}] + b_h) \\ \hat{y}^{< t>} &= g(W_{vh}h^{< t>} + b_v) \end{split}$$

where  $[h^{<l-1>}, x^{<l>}]$  means that  $h^{<l-1>}$  and  $x^{<l>}$  are concatenated together. In the next cell we provide the implementation of the forward method for a vanilla RNN.

```
In [4]: W def forward_V_RNN(inputs, weights): # Forward propagation for a a single vanilla RNN cell
x, h_t = inputs

# weights.
wh, _, _, bh, _, _ = weights

# new hidden state
h_t = np.dot(wh, np.concatenate([h_t, x])) + bh
h_t = sigmoid(h_t)

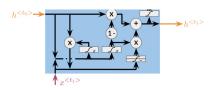
return h_t, h_t
```

As you can see, we omitted the computation of  $\hat{y}^{<\!\!\!/}$ . This was done for the sake of simplicity, so you can focus on the way that hidden states are updated here and in the GRU cell.

#### 1.2 Forward method for GRUs

A GRU cell have more computations than the ones that vanilla RNNs have. You can see this visually in the following diagram:





As you saw in the lecture videos, GRUs have relevance  $\Gamma_r$  and update  $\Gamma_u$  gates that control how the hidden state  $h^{</r>$ these gates, GRUs are capable of keeping relevant information in the hidden state even for long sequences. The equations needed for the forward method in GRUs are provided below:

$$\begin{split} \Gamma_r &= \sigma(W_r[h^{< t-1>}, x^{< t>}] + b_r) \\ \Gamma_u &= \sigma(W_u[h^{< t-1>}, x^{< t>}] + b_u) \\ c^{< t>} &= \tanh(W_h[\Gamma_r * h^{< t-1>}, x^{< t>}] + b_h) \\ h^{< t>} &= \Gamma_u * c^{< t>} + (1 - \Gamma_u) * h^{< t-1>} \end{split}$$

In the next cell, please implement the forward method for a GRU cell by computing the update u and relevance r gates, and the candidate hidden state c.

```
In [5]: 
 \mbox{\it M} def forward_GRU(inputs, weights): # Forward propagation for a single GRU cell x, h_t = inputs
                              # weights.
                              wu, wr, wc, bu, br, bc = weights
                             # Update gate
### START CODE HERE (1-2 LINES) ###
u = np.dot(wu, np.concatenate([h_t, x])) + bu
u = sigmoid(u)
                              ### END CODE HERE ###
                             ### TART CODE HERE (1-2 LINES) ###
r = np.dot(wr, np.concatenate([h_t, x])) + br
r = sigmoid(r)
                              ### END CODE HERE ###
                             # Candidate hidden state
### START CODE HERE (1-2 LINES) ###
c = np.dot(wc, np.concatenate([r * h_t, x])) + bc
c = np.tanh(c)
### END CODE HERE ###
                             # New Hidden state h_t
h_t = u* c + (1 - u)* h_t
return h_t, h_t
```

Run the following cell to check your implementation.

```
Out[6]: array([[ 9.77779014e-01], [-9.97986240e-01],
                      -5.19958083e-011.
                      -9.99999886e-01],
-9.99707004e-01],
                       -3.02197037e-041.
                       -9.58733503e-011
                       2.10804828e-02],
9.77365398e-05],
                       9.99833090e-011.
                       1.63200940e-08],
8.51874303e-01],
                       5.21399924e-021.
                       2.15495959e-02],
9.99878828e-01],
                     9.77165472e-0111)
         Expected output:
             array([[ 9.77779014e-01],
                     [-9.97986240e-01],
                     [-5.19958083e-01],
                     [-9.99999886e-01].
                     [-9.99707004e-01],
                     [-3.02197037e-04],
                     [-9.58733503e-01],
                     [ 2.10804828e-02].
                     [ 9.77365398e-05],
                     [ 9.99833090e-01],
                     [ 1.63200940e-08],
                       8.51874303e-01],
```

# Part 2: Implementation of the scan function

In the lectures you saw how the scan function is used for forward propagation in RNNs. It takes as inputs:

- fn : the function to be called recurrently (i.e. forward\_GRU )
- elems : the list of inputs for each time step (  $\ensuremath{\mathsf{X}}$  )

[ 5.21399924e-02], [ 2.15495959e-02], 9.99878828e-01], [ 9.77165472e-01]])

- · weights : the parameters needed to compute fn
- h\_0 : the initial hidden state

 $scan \ goes \ through \ all \ the \ elements \ x \ in \ elems \ , \ calls \ the \ function \ fn \ with \ arguments \ ([\ x \ , \ h_t\ ], \ weights \ ), \ stores \ the \ computed \ hidden \ state \ h_t \ )$ and appends the result to a list ys. Complete the following cell by calling fn with arguments ([ x ,  $h_t$  ], weights ).

```
In [7]: 
oldsymbol{N}
 def scan(fn, elems, weights, h_0=None): # Forward propagation for RNNs h_t = h_0 ys = [] for x in elems: ### START CODE HERE (1 LINE) ### y, h_t = fn([x, h_t], weights) ### END CODE HERE ###
```

```
ys.append(y)
return ys, h_t
```

## Part 3: Comparison between vanilla RNNs and GRUs

You have already seen how forward propagation is computed for vanilla RNNs and GRUs. As a quick recap, you need to have a forward method for the recurrent cell and a function like scan to go through all the elements from a sequence using a forward method. You saw that GRUs performed more computations than vanilla RNNs, and you can check that they have 3 times more parameters. In the next two cells, we compute forward propagation for a sequence with 256 time steps (T) for an RNN and a GRU with the same hidden state h\_t size (h\_dim =16).

As you were told in the lectures, GRUs take more time to compute (However, sometimes, although a rare occurrence, Vanilla RNNs take more time. Can you figure out what might cause this?). This means that training and prediction would take more time for a GRU than for a vanilla RNN. However, GRUs allow you to propagate relevant information even for long sequences, so when selecting an architecture for NLP you should assess the tradeoff between computational time and performance

Congratulations! Now you know how the forward method is implemented for vanilla RNNs and GRUs, and you know how the scan function provides an abstraction for forward propagation in RNNs.